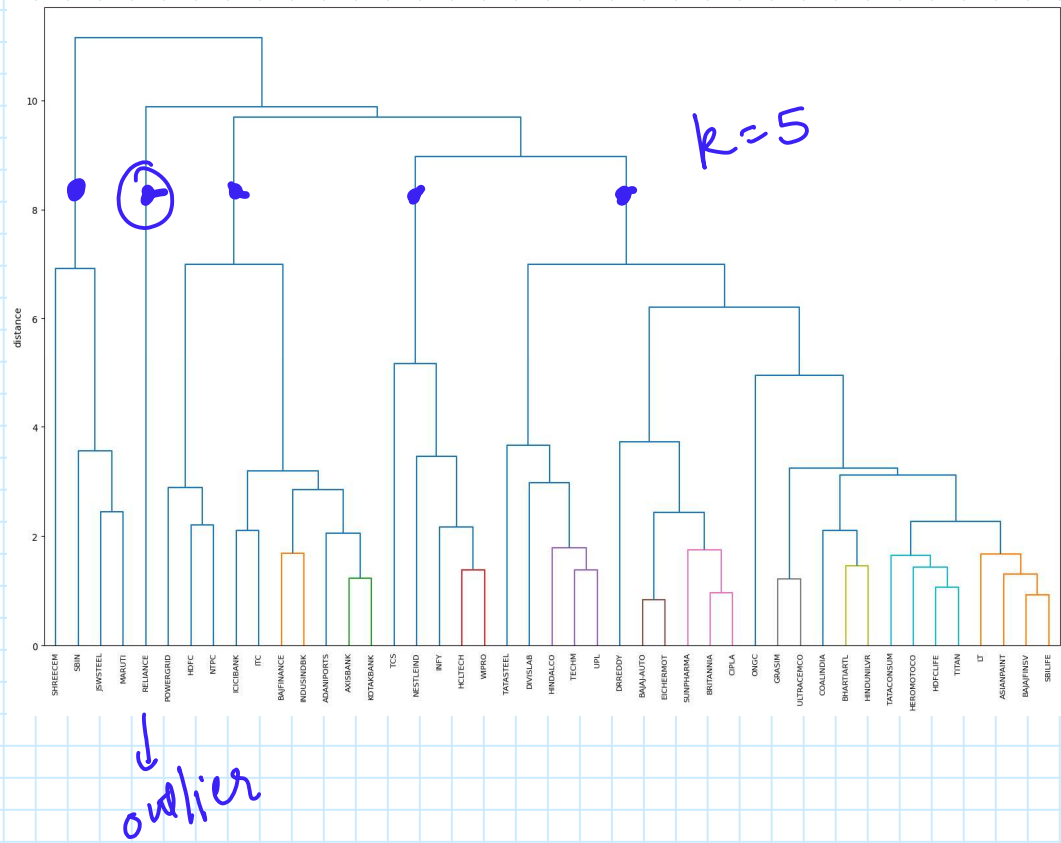
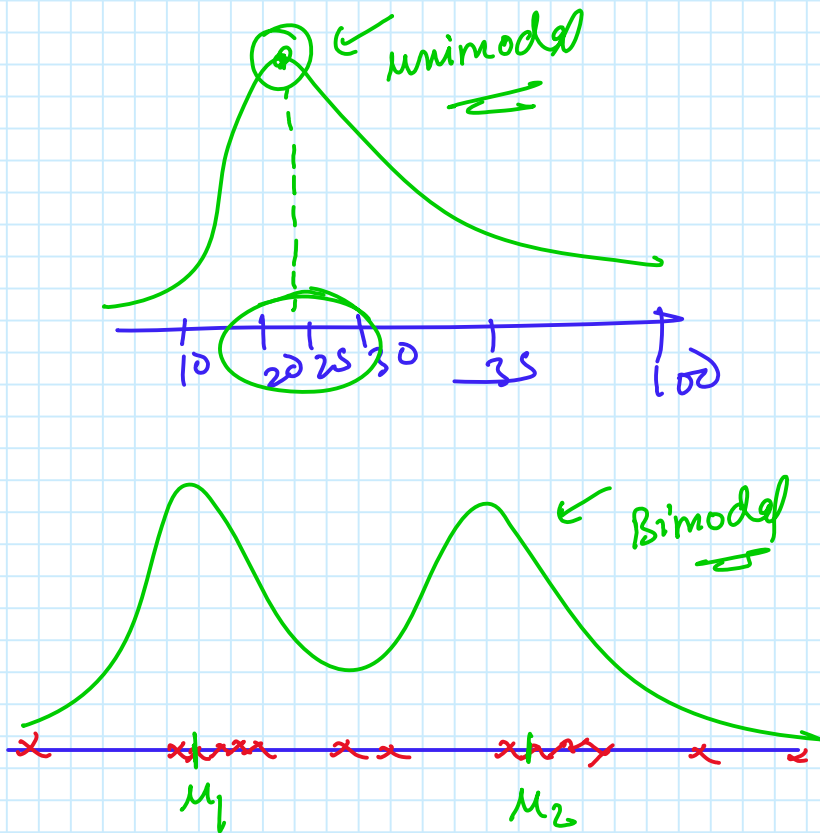
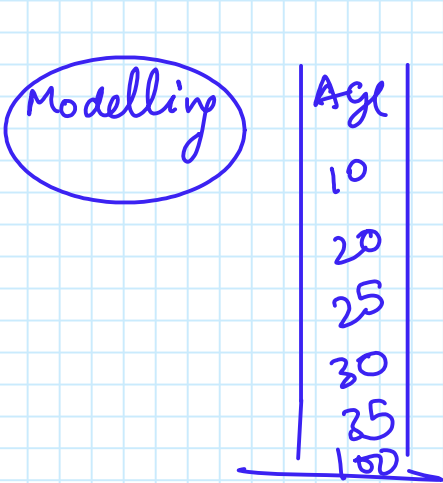
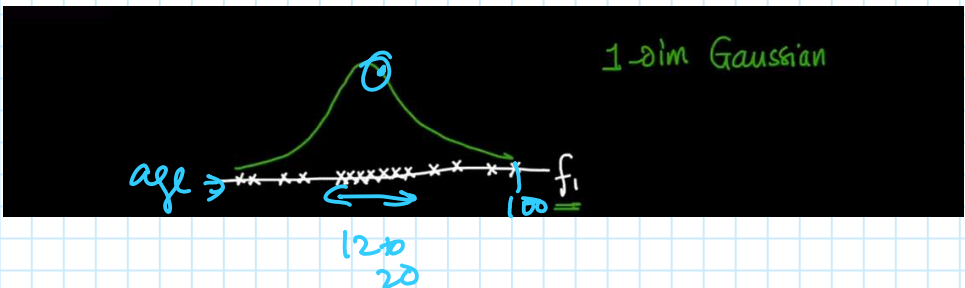


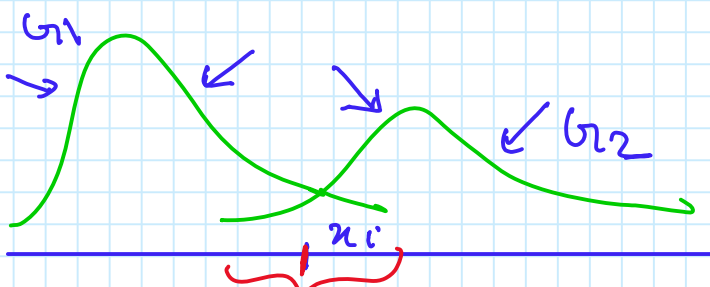
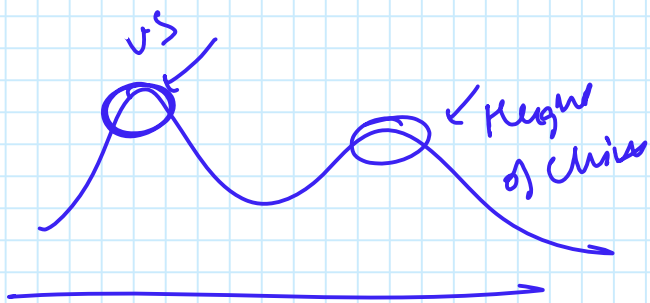
Gaussian Mixture Models



Gaussian



- In that case, there will be two **Gaussian Distributions** with mean μ_1 and μ_2 .
- If there is only a single **mount(peak)** in the distribution of a data, then the data is known as **uni-modal data**
- If there are more than one **mount(peak)** in the distribution of a data, then the data is known as **multi-modal data**



If we combine both the gaussians, the point at which both distributions intersects will have more height than the start and the end of the distribution.

This is known as **Mixture of Gaussians**

Q. What does the intersection represents?

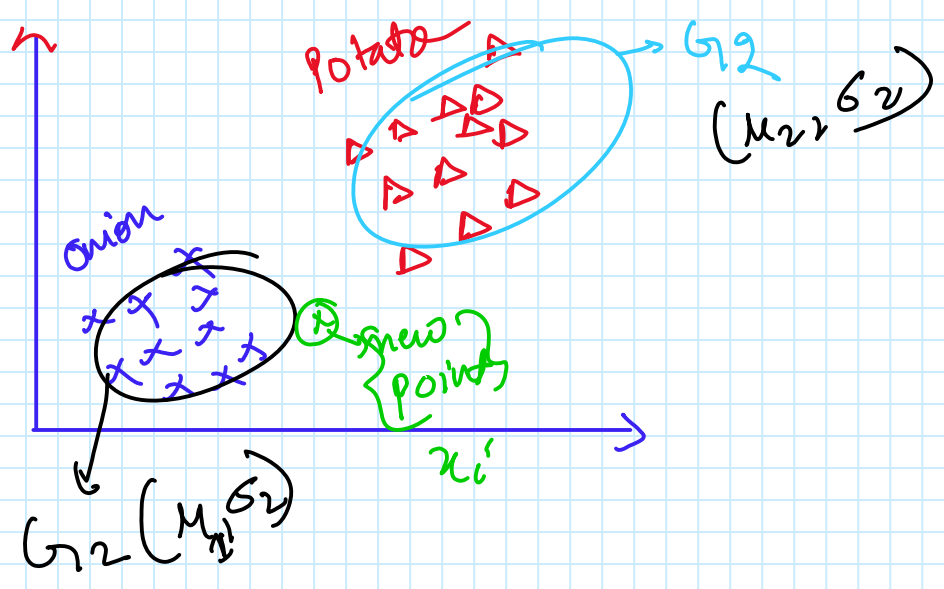
- It says that the point at intersection can either belong to distribution G_1 or distribution G_2

$P(G_1|x_i) = 0.02$ and $P(G_2|x_i) = 0.06$

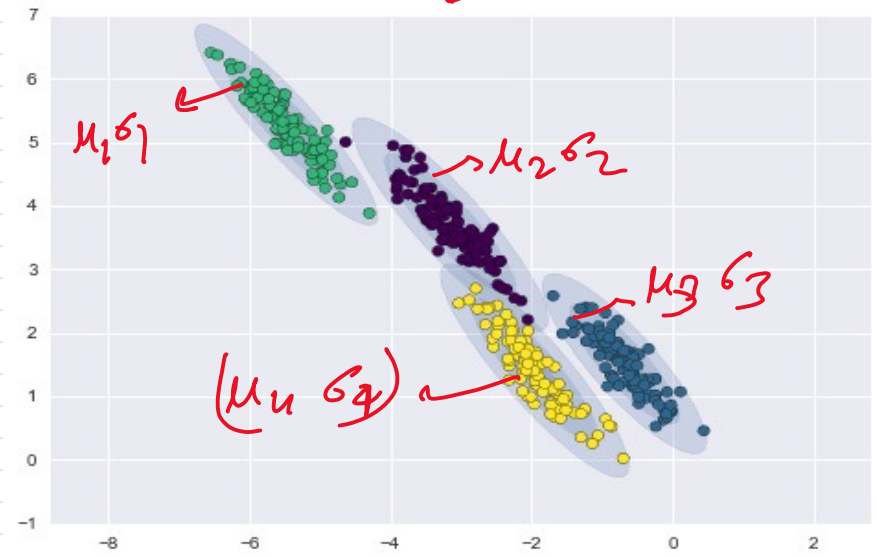
Normalise $\Rightarrow \frac{0.02}{0.02+0.06} = 20\%$ $\frac{0.06}{0.02+0.06} = 80\%$



Gaussian Mixture Models → model data as mixture of gaussians



$k=4$



$k \rightarrow$ means \rightarrow Hard Clustering
Onion \rightarrow Potato

$$P(x_i = G_1) = \frac{1}{\sqrt{2\pi}\sigma_1} e^{-\frac{(x-\mu_1)^2}{2\sigma_1^2}}$$

$$P(x_i = G_2) = \frac{1}{\sqrt{2\pi}\sigma_2} e^{-\frac{(x-\mu_2)^2}{2\sigma_2^2}}$$

$\Rightarrow 0.03$

Normalised

$$\frac{0.03}{\text{Total} = (0.03 + 0.4)} = 3\%$$

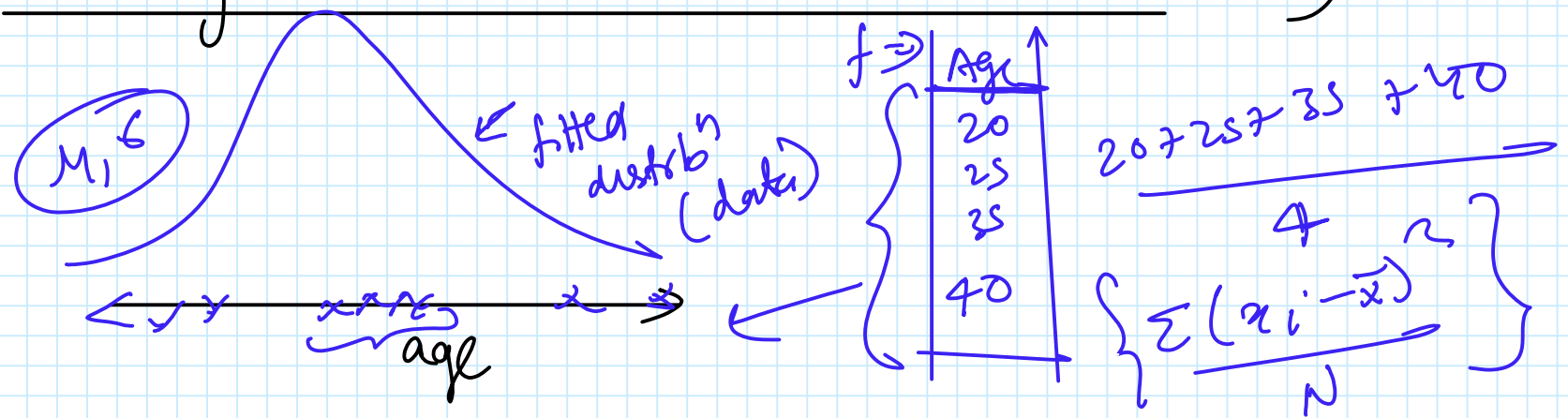
$\Rightarrow 0.4$

$$\frac{0.4}{\text{Total} (0.03 + 0.4)} = 97\%$$

Gaussian \rightarrow (completely characterised by Mean & Std Dev)

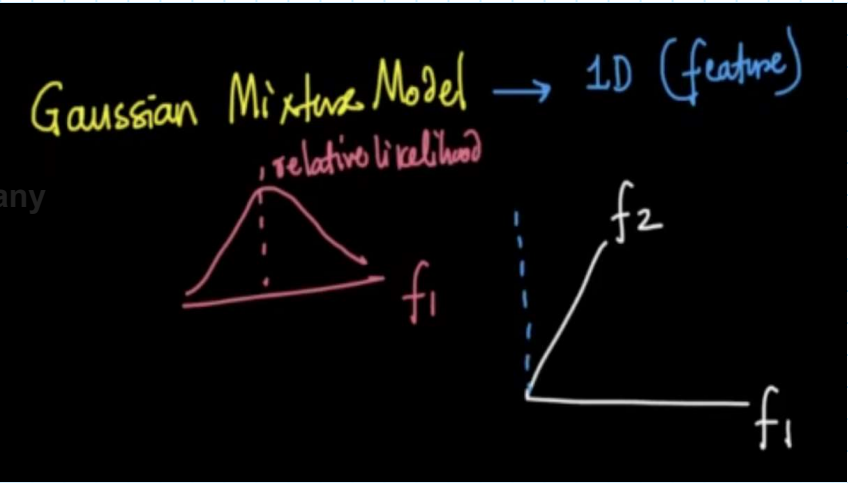


Extending Gaussian to multi Dimensions (2D)



Intuition behind Gaussian Mixture Model

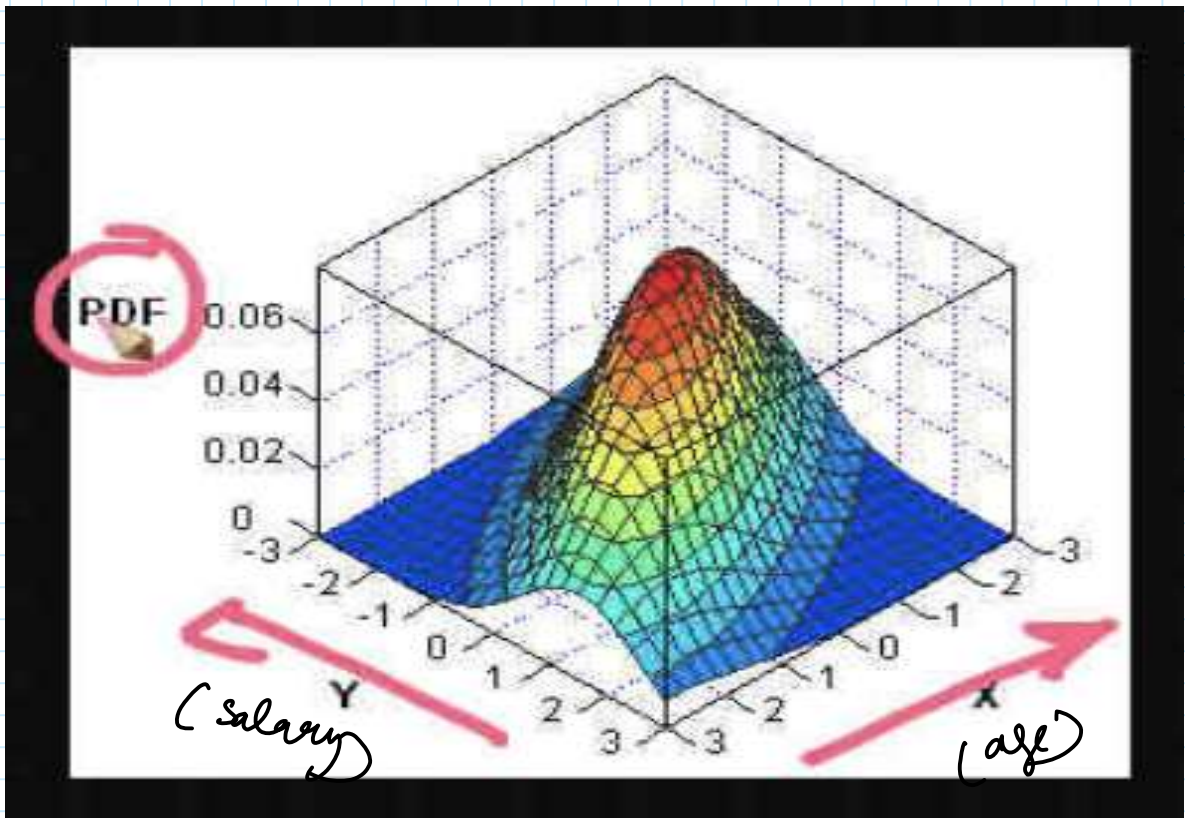
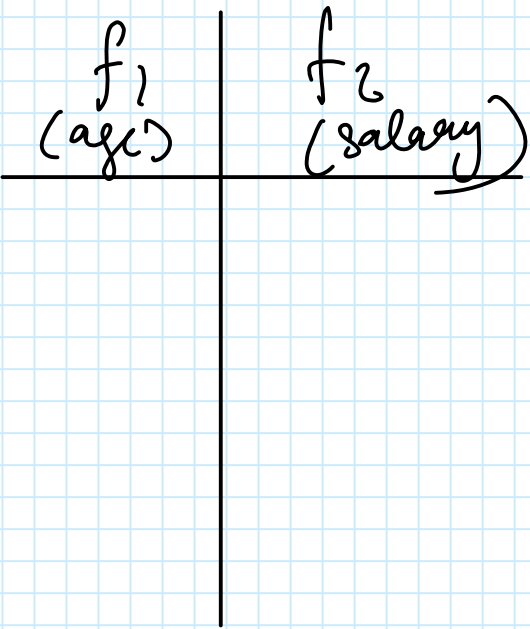
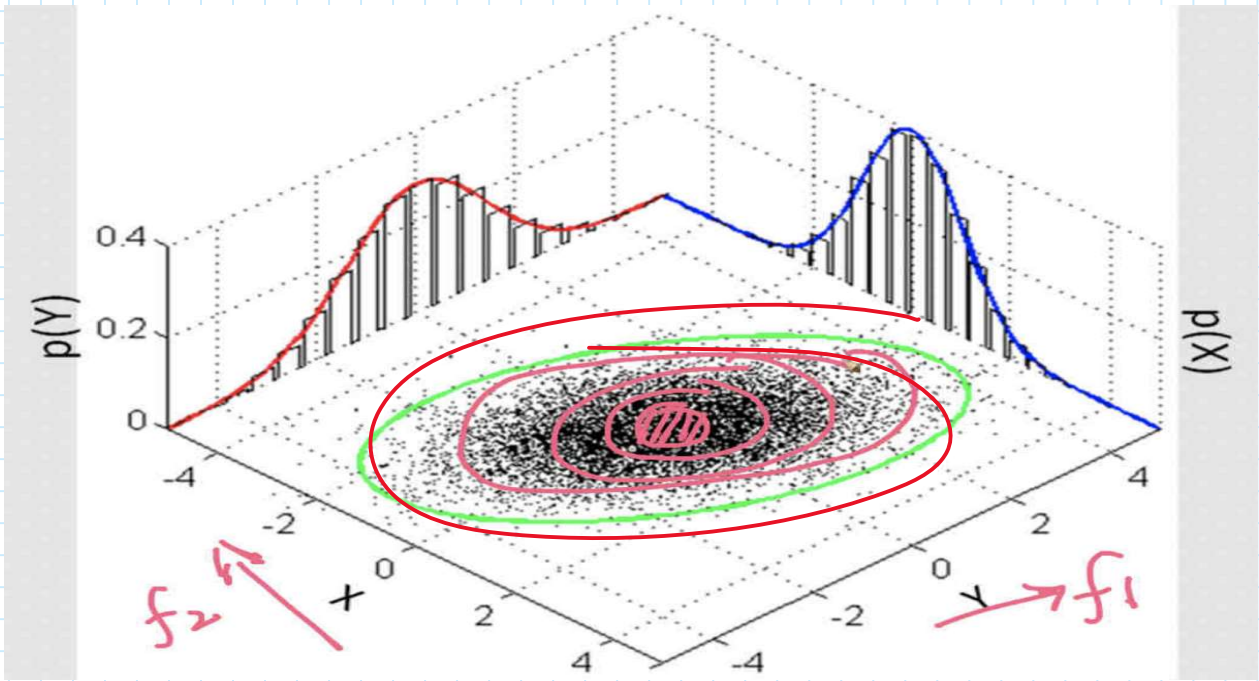
- The basic intuition behind Gaussian Mixture Model is that we can model any data, that we have, as a mixture of Gaussians.



GAUSSIAN IN 2 Dimensions

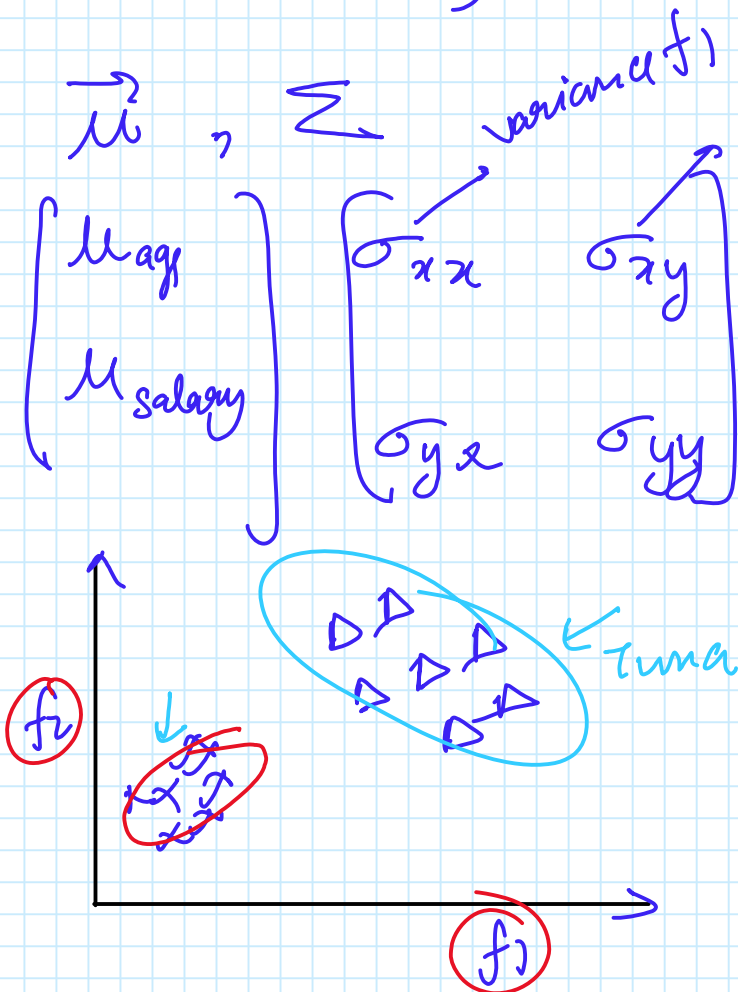
- As we can see, the data points are more dense in the centre and as we move away from the centre, the density of points decreases. This is to simulate a Gaussian (Normal) Distribution.
- If we consider only 1 feature at a time, we can see its Gaussian Distribution:
 - Red curve represents the Gaussian Distribution for Feature 1 alone.
 - Blue curve represents the Gaussian Distribution for Feature 2 alone.
- The Gaussian Distribution of the 2 features combined would look something like shown below:

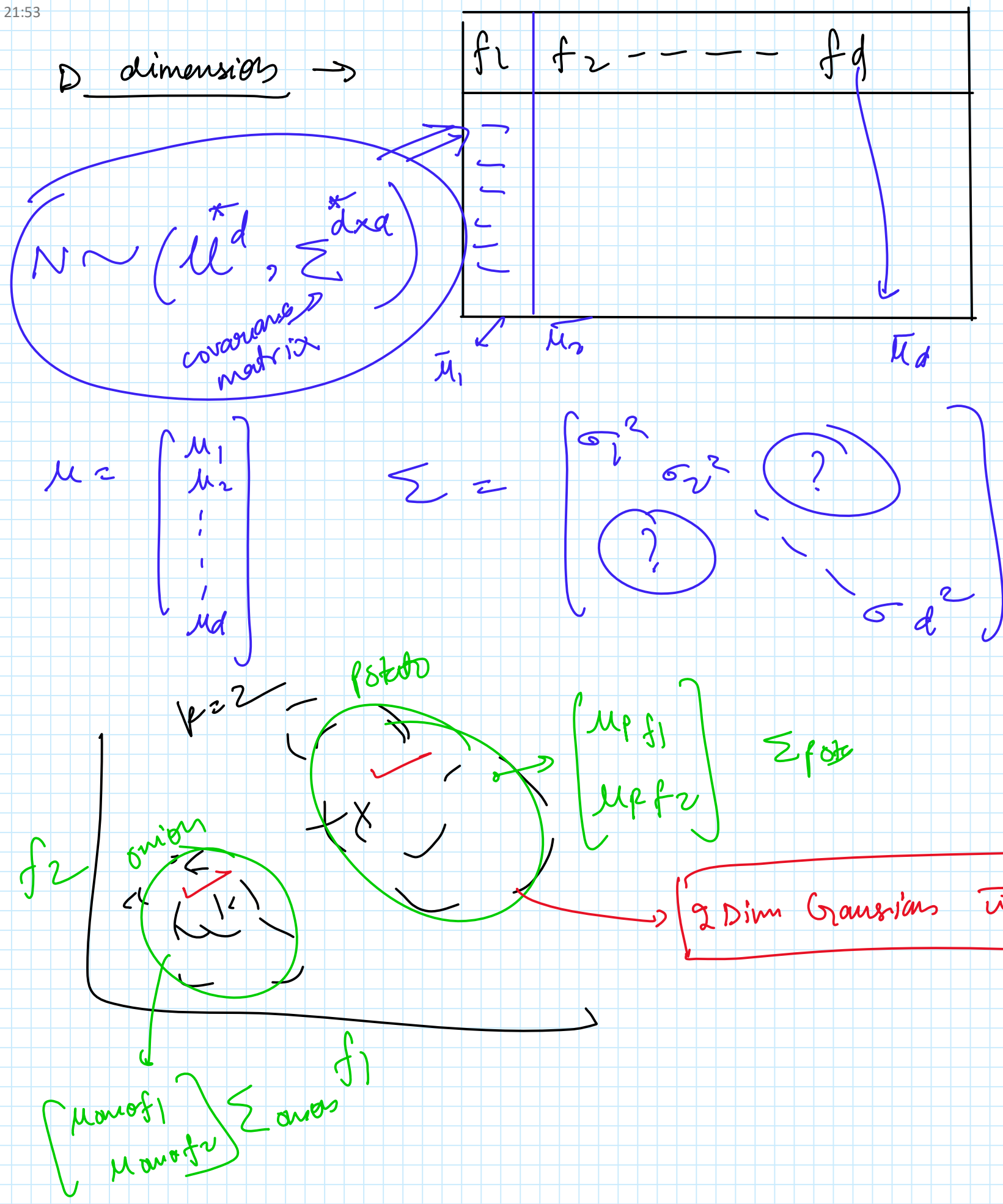
* [Gaussian Distibⁿ for Two features]



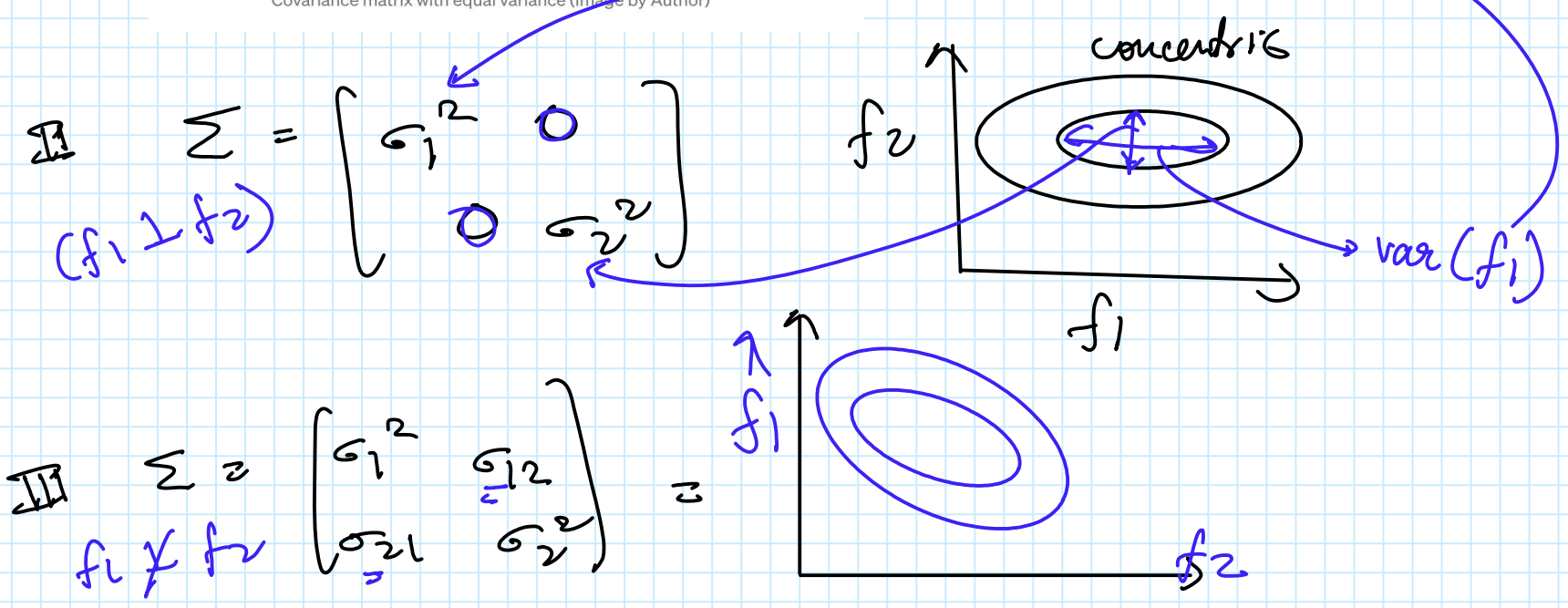
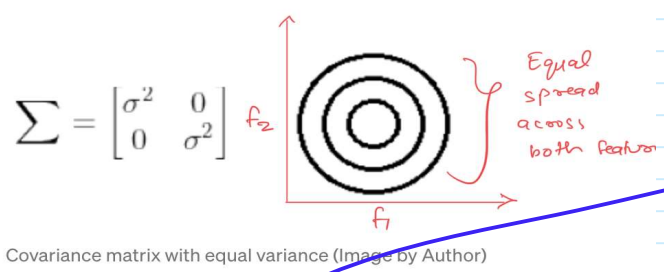
μ, σ \rightarrow μ potaka μ aniva Σ

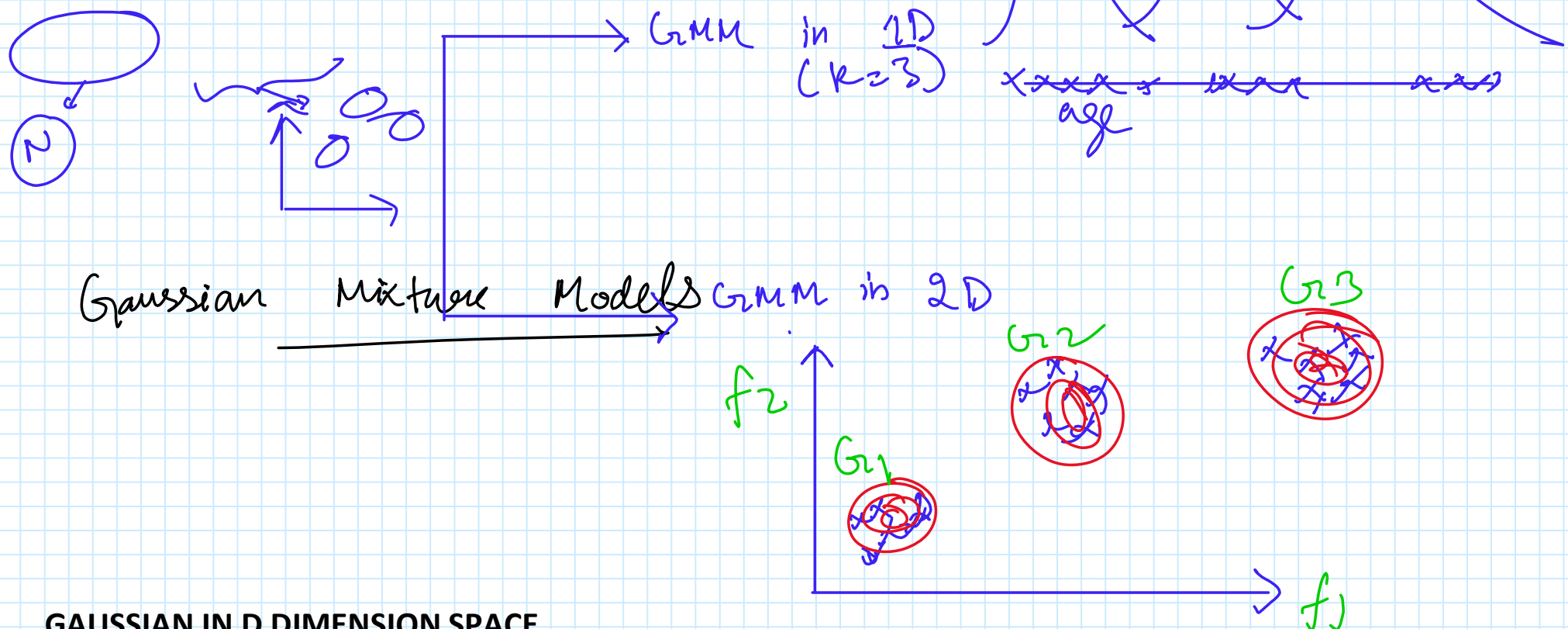
Parameters (2D)





Ques

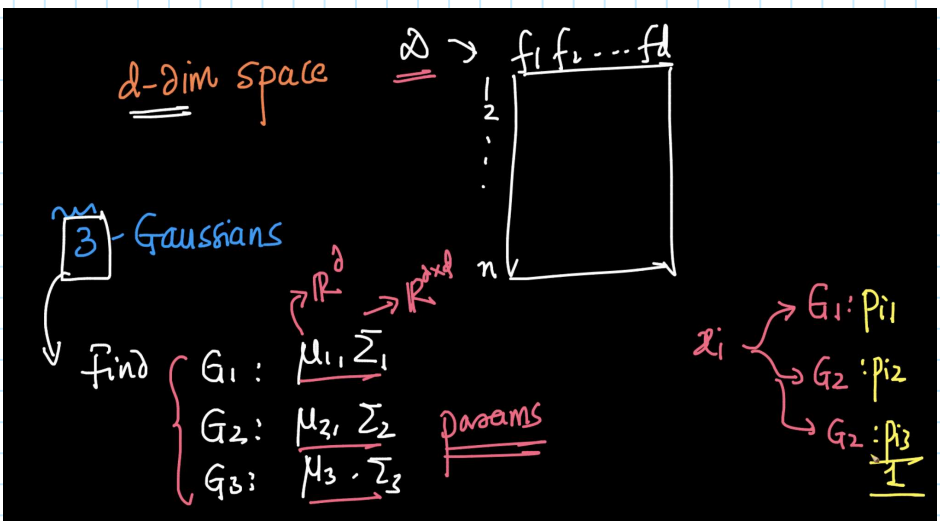




GAUSSIAN IN D DIMENSION SPACE

(D → dimensional space)
Each gaussian will have its own set of parameters

- $G_1 : \mu_1, \Sigma_1$
- $G_2 : \mu_2, \Sigma_2$
- $G_3 : \mu_3, \Sigma_3$



x_i → $G_1 (0.2)$ → 0.2 / 0.8 %
 x_i → $G_2 (0.1)$ → 0.1 / 0.8 %
 x_i → $G_3 (0.5)$ → 0.5 / 0.8 %

- For a d-Dimensional Gaussian Mixture Model, we'll need:
- The d-dimensional vector of means $\mu(d)$ and the $d \times d$ Covariance Matrix $\Sigma(d \times d)$, i.e., the Normal Distribution N_d .
 - If we have this Normal Distribution N_d , we can compute the probability of a data point belonging to a Gaussian (Cluster) using a slightly more complex density function:

$P(x_i \rightarrow G_1) =$ $\mu_d, \Sigma_{d \times d}$ \times (Ignore)

definite. In this case the distribution has density^[5]

$$f_{\mathbf{x}}(x_1, \dots, x_d) = \frac{\exp(-\frac{1}{2}(\mathbf{x} - \underline{\mu})^T \Sigma^{-1}(\mathbf{x} - \underline{\mu}))}{\sqrt{(2\pi)^d |\Sigma|}}$$

$\sqrt{(2\pi)^d |\Sigma|}$ → determinant

1D formula → $\frac{1}{\sqrt{2\pi}\sigma^2} e^{-\frac{1}{2}(\frac{x-\mu}{\sigma})^2}$

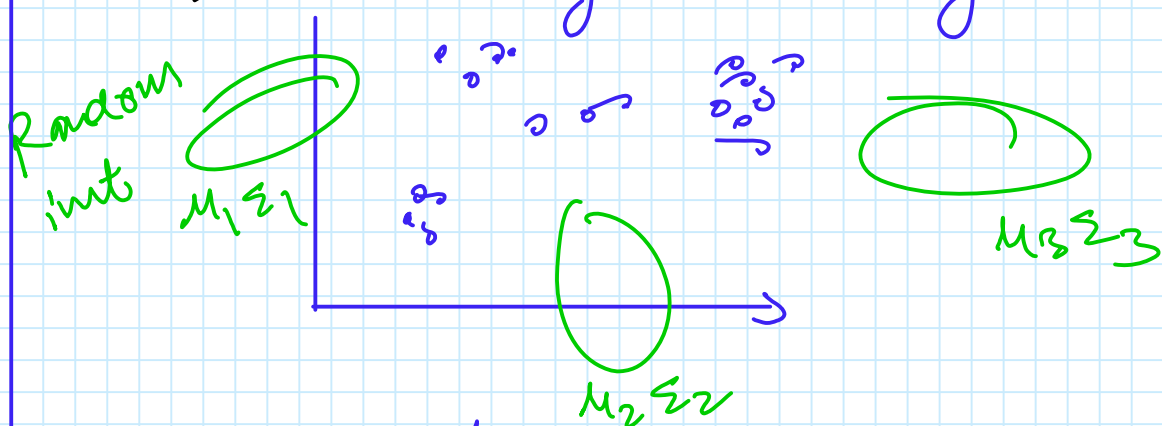
CPD f of

EM find the parameters of model

Expectation - maximisation algorithm

$(p \sim i)$

$E-M \rightarrow$ Randomly initialize gaussians (μ, Σ) for each



Step 2 \rightarrow Assign/find prob of each pt belonging to each gaussian

→ for each $x_i \rightarrow 1$ to N

$$x_i \rightarrow \left\{ \begin{array}{l} G_1 \rightarrow 0.2 \\ G_2 \rightarrow 0.3 \\ G_3 \rightarrow 0.4 \end{array} \right\}$$

Diagram illustrating a branching process starting from a root node x_1 . Three arrows branch out to nodes g_1 (20%), g_2 (30%), and g_3 (50%). A dashed line leads down to a node x_N , which is connected to the root by a large curved arrow on the left.

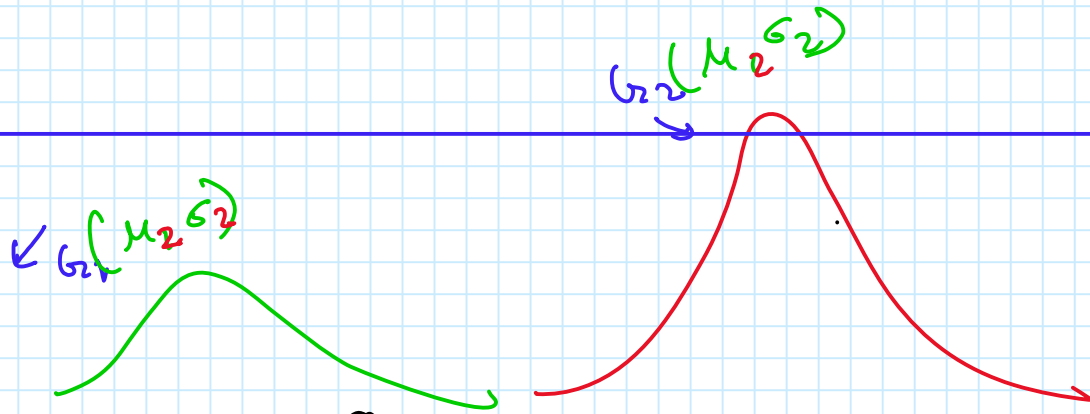
Step 3

Update the centroid
(K Means)

updates the gaussians $\rightarrow \pi$
(GMM)

$$\vec{\mu}$$

GMM Visualise



Agg
Agg
(P=22) ASS

Assign →

update

formulas

k mean.

$$\frac{\sum x_i}{N} \quad i \in C_1$$

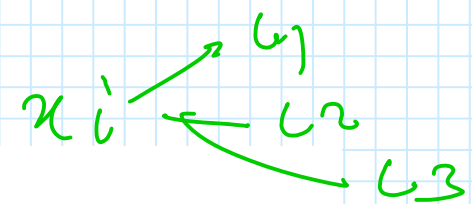
update formula \rightarrow weighted Average $\rightarrow \frac{\sum w_i x_i}{N} / \frac{\sum (p_i(x_i))}{N} \times \left(\frac{\sum x_i}{N} \right)$

→ K means

(E, M) →
Random
init gaussians

Step-1 Expectation:

- This is basically an assignment step.
- In this step, for each point x_i , we compute the probability of point x_i belonging to a j^{th} cluster
- So, initially we start by randomly assigning a probability such that it belongs to j^{th} gaussian (j^{th} cluster).
- We then compute the probabilities of a point belonging to k different clusters using Probability Density Function



Step-2 Maximization:

- In GMMs, we compute normalized probabilities of each and every point belonging to a cluster.
- This is kind of like a soft assignment, whereas in KMeans we did the Hard Assignment by randomly picking centroids.
- In the Maximization step we re-estimate gaussian parameters, for all K gaussians.

$$\mu_j' = \frac{\sum p_i x_i}{N}$$

K Means → $\frac{\sum x_i}{N}$ → In GMM → $\frac{\sum p_i x_i}{N}$

⊗ Very high Dimensional → Imbalance issue

GMM Algorithm & Implementation

There are two steps involved in GMM Algorithm:

1. Expectation
2. Maximization

Step-1. Expectation:

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- In this step, for each point x_i , we compute the probability of point x_i belonging to a j^{th} cluster
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Q. How do we do that?

- We update mean vector (μ_j) and covariance vectors (Σ_j) using weighted scheme

What is weighted scheme?

- Suppose we've to update mean vector (μ_j)
- Instead of taking simple average as we do in K-Means, we compute weighted average of each x_i , where the weightage is based on the probabilities that the point $x_i \in$ cluster j
- We also compute covariance vector in the same fashion
- In the second step of K-means, we update the points based on the updated centroids of the cluster
- Whereas in GMM, we update a point based on the mean (which is same as centroid in K-Means) and covariance. The only difference is that we use weighting scheme in GMM Models